In the Claims:

Claim 1 (Currently amended): A compound of the formula:

in which:

- a) X¹ is represented by cyano, halogen or haloalkyl,
- b) one of R^1 or R^2 is represented by C_1 - C_6 alkyl which may be optionally substituted, and the other of R^1 or R^2 is represented by hydrogen or C_1 - C_6 alkyl which may be optionally substituted.
- c) Alk¹ is represented by a C₁-C₂ linear alkylene group, in which up to two hydrogen atoms are optionally replaced by a substituent selected from the group consisting of C₁-C₆ alkyl optionally substituted, halogen, hydroxy, thiol, and cyano,
- d) n is represented by the integer 0 or 1,
- e) Y is represented by NX²X³-or-O-X³,
- f) X^2 is represented by hydrogen or $(C_{1}-C_{6})$ alkyl optionally substituted, <u>and</u>.
- g) X³ is represented by

i.hydrogen.

 $\mathsf{ii.}(\mathsf{C_{4-}C_{42-}}) \mathsf{alkyl}, \ \mathsf{optionally\ substituted},$

iii. $(C_{_2}$ - $C_{_{12}}$)alkenyl, optionally substituted,

iv. $(C_2 - C_{42})$ alkynyl, optionally substituted,

v. (C₃-C₁₀)cycloalkyl, optionally substituted,

vi. (C_3-C_{10}) cycloalkyl (C_4-C_6) alkyl, in which the alkyl and cycloalkyl moieties may each be optionally substituted,

 $\underline{\text{vii.i.}} \quad \text{(C}_{\scriptscriptstyle 6}\text{-C}_{\scriptscriptstyle 10}\text{)aryl, optionally substituted,}$

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- - fii.-(CH₂)-(Alk²)_q-C(O)R³, in which Alk² is represented by a (C₁-C₈) linear alkylene group, in which up to eight hydrogen atoms may optionally be replaced by a substituent, selected from the group consisting of (C₁-C₆) alkyl optionally substituted, (C₁-C₆) alkoxy, halogen, hydroxy, thiol, cyano, and NR⁸R⁹ in which R⁸ and R⁹ are each independently represented by hydrogen or (C₁-C₆) alkyl, q is the integer 0 or 1, R³ is represented by hydrogen, (C₄-C₁₂)alkyl, (C₆-C₁₆)aryl, or (C₆-C₁₀)aryl(C₄-C₆)alkyl, in which the alkyl and aryl moieties may each be optionally substituted,
 - iv. (CH_2) - $(Alk^2)_q$ -C(O)-O- R^4 , in which Alk^2 and q_r are as defined above, and R^4 is represented by hydrogen, $(C_4$ - C_{42})alkyl, $(C_6$ - C_{40})aryl, or $(C_6$ - C_{40})aryl $(C_4$ - C_6)alkyl, in which the alkyl and aryl moieties may be optionally substituted.
 - v:-(CH₂)-(Alk²)_q-C(O)-NR⁵R⁶ in-which Alk² and q are as described above, and R⁵ and R⁶ are each independently represented by hydrogen, (C₁-C₁₂)alkyl, (C₆-C₄₀)aryl, or (C₆-C₄₀)aryl(C₁-C₆)alkyl, in which the alkyl and aryl moieties may be optionally substituted.
 - vi.- (CH_2) - $(Alk^2)_q$ -Y- R^7 , in which Alk^2 and q are as defined above, Y is O or S, and R^7 is selected from the group consisting of hydrogen, $(C_4$ - C_{42})alkyl, $(C_6$ - C_{10})aryl, or $(C_6$ - C_{40})aryl $(C_4$ - C_6)alkyl, in which the alkyl and aryl moieties may be optionally substituted.

vii.heteroaryl, optionally substituted,

viii.heteroaryl(C₁_C₆)alkyl, in which the heteroaryl and alkyl moieties may each be optionally substituted.

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xv.iii. _-heterocyclic, optionally substituted,
x.heterocyclic(C₄_C₆)alkyl, in which the alkyl
and heterocyclic moieties may each be substituted, or,

h) for those compounds in which Y is N, X² and X³, along with the adjacent nitrogen atom, may form a heterocyclic ring, which may optionally be substituted, or a pharmaceutically acceptable salt, or solvate, thereof.

Claim 2 (Original): A compound according to claim 1 in which one of R^1 or R^2 is hydrogen and the other of R^1 or R^2 is selected from the group consisting of isobutyl, propyl, n-butyl, isopropyl, and ethyl.

Claim 3 (Previously amended): A compound according to claim 2 in which n is 0.

Claim 4 (Currently amended): A compound according to claim 3 in which X¹ is trifluoromethyl and is located at the 3-position of the phenyl ring.

Claim 5 (Cancelled)

Claim 6 (Currently amended): A compound according to claim 4 in which χ^2 is hydrogen.

Claim 7 (Cancelled)

Claim 8 (Cancelled)

Claim 9 (Currently amended): A compound according to anyone of claim 18 in which X¹ is represented by halogen or haloalkyl.

Claim 10-12 (Cancelled)

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Claim 13 (Previously amended): A pharmaceutical composition comprising a compound according to claim 1 in admixture with 1, or more, pharmaceutically acceptable excipients.

Claim 14 (Previously amended): A topical pharmaceutical formulation comprising a compound according to claim 1 in admixture with 1, or more, pharmaceutically acceptable excipients suitable for dermal application.

Claim 15 (Previously amended): A kit comprising a compound according to claim 1 packaged for retail distribution, which advises a consumer how to utilize the compound to alleviate a condition selected from the group consisting of acne, alopecia, and oily skin.

Claim 16. (New): A compound according to claim 1 in which X¹ is represented by CF₃ and is located at the 3-position of the phenyl ring, R¹ is isobutyl or propyl, R² is hydrogen, and n is 0.

Claim 17. (New): A compound according to claim 1 in which X^1 is represented by CF_3 and is located at the 3-position of the phenyl ring, R^1 is isobutyl, R^2 is hydrogen, n is 0, and X^2 is represented by hydrogen.

Claim 18 (New): A compound according to claim 1 in which X^1 is represented by CF_3 and is located at the 3-position of the phenyl ring, R^1 is isobutyl or propyl, R^2 is hydrogen, n is 0, X^2 is represented by hydrogen and X^3 is benzyl or phenethyl in which the phenyl ring is optionally substituted with at least one substituent selected from the group consisting of methoxy, ethoxy, hydroxy, and methyl.

Claim 19 (New): A compound according to claim 1 selected from the group consisting of:

- a) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid benzylamide,
- b) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid benzylamide,

- c) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid [2-(4-methoxy-phenyl)-ethyl]-amide,
- d) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid (2-phenoxy-ethyl)-amide,
- e) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid -3-methoxy-benzylamide,
- f) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid [2-(4-methoxy-phenyl)-ethyl]amide,
- g) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-2-methoxy-benzylamide,
- h) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-2-ethoxy-benzylamide,
- i) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-3-methylbenzylamide, and,
- j) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-2-methylbenzylamide.

Claim 20 (New): A compound according to claim 1 selected from the group consisting of:

- a) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-4-methoxy-benzylamide,
- b) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-3-methoxy-benzylamide,
- c) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-2-methoxy-benzylamide,
- d) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-2-ethoxy-benzylamide,
- e) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-3-methyl-benzylamide,
- f) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-2-methyl-benzylamide,
- g) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-2,4-dimethyl-benzylamide,
- h) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-4-methoxy-benzylamide,
- i) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-(2-p-tolyl-ethyl)-amide, and,
- j) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-[2-(2-methoxy-phenyl)-ethyl]-amide.

Claim 21 (New): A compound according to claim 1 selected from the group consisting of:

- a) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-(2-m-tolylethyl)-amide,
- b) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-(2-p-tolyl-ethyl)-amide,
- c) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-[2-(2-methoxy-phenyl)-ethyl]-amide,
- d) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-(2-m-tolyl-ethyl)-amide,
- e) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-(2-phenoxy-propyl)-amide,
- f) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-(2-phenoxy-ethyl)-amide.
- g) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-[2-(3-methoxy-phenyl)-ethyl]-amide,
- h) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-[2-(4-hydroxy-phenyl)-ethyl]-amide,
- i) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid benzyl-isopropyl-amide,
- j) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-[2-(3-methoxy-phenyl)ethyl]-amide,
- k) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-[2-(4-hydroxy-phenyl)-ethyl]-amide,
- 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid benzyl-isopropylamide.
- m) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid [1-(4-hydroxy-phenyl)-ethyl]-amide,
- n) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid 4-isopropyl-benzylamide,
- o) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-3-methoxy-benzylamide, and,
- p) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-4-methoxy-benzylamide.

Claim 22 (New): A compound according to claim 1 selected from the group consisting of:

- a) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-3,4-dihydroxy-benzylamide,
- b) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid (naphthalene-1-yl-methyl)-amide,
- c) 2-(4-cyano-3-trifluoromethyl-phenoxy)-hexanoic acid benzylamide,
- d) N-benzyl-2-(4-cyano-3-trifluoromethyl-phenoxy)-3-methyl-butyramide,
- e) (R)- 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid benzylamide,
- f) (R)-2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid benzylamide,
- g) (R)-2-(cyano-3-trifluoromethyl-phenoxy)-pentanoic acid 2-methylbenzylamide,
- h) 2-(3-chloro-4-cyano-phenoxy)-pentanoic acid benzylamide, and,
- 2-(3-chloro-4-cyano-phenoxy)-pentanoic acid [2-(4-hydroxy-phenyl)-ethyl]amide.

Claim 23 (New): A compound according to claim 1 selected from the group consisting of:

- a) (S)-2-(3-chloro-4-cyano-phenoxy)-pentanoic acid benzylamide,
- b) (\$)-2-(3-chloro-4-cyano-phenoxy)-pentanoic acid [2-(4-hydroxy-phenyl)ethyl]amide,
- c) (S)-2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid-3-methyl-benzylamide,
- d) (S)-2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-[2-(4-hydroxy-phenyl)-ethyl]-amide,
- e) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid-[1-(methoxy-phenyl)-ethyl]-amide,
- f) (R-) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid [2-4-hydroxy-phenyl)-ethyl]-amide,
- g) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-penanoic acid (1-phenyl-ethyl)-amide,

- h) 2-(4-cyano-3-trifluoromethyl-phenoxy)-penanoic acid[1-(4-methoxy-phenyl)-ethyl]-amide,
- i) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid [1-(4-hydroxy phenyl)-ethyl]-amide,
- j) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid [1-(4-hydroxy-phenyl)-ethyl]-amide,
- k) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid [1-(4-hydroxy-[phenyl)-ethyl]-amide,
- I) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid [1-hydroxy-phenyl)-ethyl]-amide,
- m) 2-(4-cyano-3-trifluoromethyl-phenoxy)-4-methyl-pentanoic acid (1-phenyl-ethyl)-amide, and,
- n) 2-(4-cyano-3-trifluoromethyl-phenoxy)-pentanoic acid [1-(methoxy-phenyl)-ethyl]-amide.